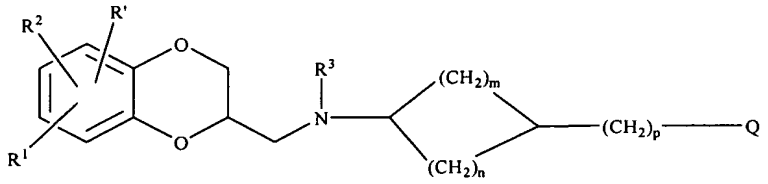
This listing of claims will replace all prior versions, and listings, of claims in the application.

Listing of Claims

1. (***currently amended***) A compound of Formula I:



I

wherein

~~**R',**~~ R^1 and R^2 are, independently, hydrogen, halo, cyano, carboxamido, carboalkoxy of two to six carbon atoms, trifluoromethyl, alkyl of 1 to 6 carbon atoms, alkoxy of 1 to 6 carbon atoms, alkanoyloxy of 2 to 6 carbon atoms, amino, mono- or di-alkylamino in which each alkyl group has 1 to 6 carbon atoms, alkanamido of 2 to 6 carbon atoms, or alkanesulfonamido of 1 to 6 carbon atoms; or R^1 and R^2, taken together, form methylenedioxy, ethylenedioxy or propylenedioxy;

R' is halo, cyano, carboxamido, carboalkoxy of two to six carbon atoms, trifluoromethyl, alkyl of 1 to 6 carbon atoms, alkoxy of 1 to 6 carbon atoms, alkanoyloxy of 2 to 6 carbon atoms, amino, mono- or di-alkylamino in which each alkyl group has 1 to 6 carbon atoms, alkanamido of 2 to 6 carbon atoms, or alkanesulfonamido of 1 to 6 carbon atoms;

R^3 is hydrogen or alkyl of 1 to 6 carbon atoms;

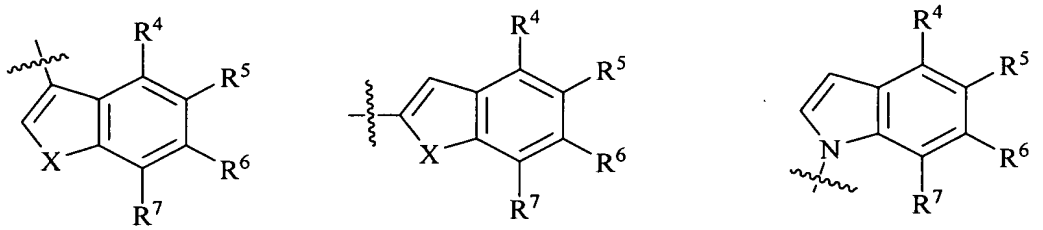
m is 1 to 3;

n is 1 or 2;

p is 0 to 3

with the proviso that when p is 0, both m and n may not be 2;

Q is a heteroaryl moiety chosen from:



R^4, R^5, R^6 and R^7 are independently selected from hydrogen, halo, cyano, trifluoromethyl, alkyl of 1 to 6 carbon atoms, alkoxy of 1 to 6 carbon atoms, alkanoyloxy of 2 to 6 carbon atoms;

X is NR^8~~, O or S~~; and

R^8 is hydrogen or alkyl of 1 to 6 carbon atoms;

or a pharmaceutically acceptable salt thereof.

2. (***original***) A compound according to claim 1, wherein R^1 and R^2 are, independently, hydrogen, halo, cyano, carboxamido, trifluoromethyl, amino, alkyl of one to six carbon atoms or alkoxy of one to six carbon atoms.

3. (***original***) A compound according to claim 1, wherein R^1 is alkoxy of one to six carbon atoms and is attached to position 8 of the benzodioxan moiety.

4. (***original***) A compound according to claim 1, wherein R^2 is hydrogen.

5. (***original***) A compound according to claim 1, wherein R^4, R^5, R^6 and R^7 are independently hydrogen, halo, cyano, alkyl of one to six carbon atoms, or alkoxy of one to six carbon atoms.

6. (***original***) A compound according to claim 1, wherein R^4, R^5, R^6 and R^7 are independently hydrogen, halo or cyano.

7. (***original***) A compound according to claim 1, wherein m and n are, independently 1 or 2.

8. (***original***) A compound according to claim 1, wherein m is 1 and n is 2.

9. (***original***) A compound according to claim 1, wherein p is 0 or 1.

10. (***original***) A compound according to claim 1, wherein p is 0.

11. (***currently amended)*** A compound according to claim 1, wherein ~~X is NR^8 and~~ R^8 hydrogen or alkyl of 1 to 3 carbons.

12. (***original***) A compound according to claim 1, wherein R^3 is hydrogen or alkyl of 1 to 3 carbons.

13. (***previously presented)*** A compound according to claim 1, wherein said compound is N-[(cis)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.

14. (***original***) A compound according to claim 1, wherein said compound is N-[(1R,3S)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.

15. (***original***) A compound according to claim 1, wherein said compound is N-[(1S,3R)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.

16. (***original***) A compound according to claim 1, wherein said compound is N-[(trans)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof

17. (***original***) A compound according to claim 1, wherein said compound is N-[(1S,3S)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.

18. (***original***) A compound according to claim 1, wherein said compound is N-[(1R,3R)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.

19. (***original***) A compound according to claim 1, wherein said compound is N-[(cis)-3-(5-fluoro-1-methyl-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.

20. (***original***) A compound according to claim 1, wherein said compound is N-[(trans)-3-(5-fluoro-1-methyl-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.

21. (***original***) A compound according to claim 1, wherein said compound is 3-[(cis)-3-({[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amino)cyclopentyl]-1H-indole-5-carbonitrile or a pharmaceutically acceptable salt thereof.

22. (***original***) A compound according to claim 1, wherein said compound is 3-[(trans)-3-({[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amino)cyclopentyl]-1H-indole-5-carbonitrile or a pharmaceutically acceptable salt thereof.

23. (***original***) A compound according to claim 1, wherein said compound is 3-[(cis)-3-({[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amino)cyclopentyl]-1-methyl-1H-indole-5-carbonitrile or a pharmaceutically acceptable salt thereof.

24. (***original***) A compound according to claim 1, wherein said compound is 3-[(trans)-3-({[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amino)cyclopentyl]-1-methyl-1H-indole-5-carbonitrile or a pharmaceutically acceptable salt thereof.

25. (***original***) A compound according to claim 1, wherein said compound is N-{[(2S)-8-ethoxy-2,3-dihydro-1,4-benzo-dioxin-2-yl]methyl}-N-[(cis)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]amine or a pharmaceutically acceptable salt thereof.

26. (***original***) A compound according to claim 1, wherein said compound is N-{[(2S)-8-ethoxy-2,3-dihydro-1,4-benzo-dioxin-2-yl]methyl}-N-[(trans)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]amine or a pharmaceutically acceptable salt thereof.

27. (***cancelled***)

28. (***cancelled***)

29. (***cancelled)***

30. (***cancelled***)

31. (***original***) A pharmaceutical composition, comprising:

an effective amount of a compound according to claim 1 or a pharmaceutically acceptable salt thereof; and

a pharmaceutically acceptable carrier or excipient.